

3,7,12-Trihydroxy-5-cholenoic acid, methyl ester, TMS

Inchi: InChI=1S/C34H64O5Si3/c1-23(14-17-31(35)36-4)26-15-16-27-32-28(22-30(34(26,27)3)3
InchiKey: VIATWLYTLFTFJF-OWWARJOYSA-N
Formula: C34H64O5Si3
SMILES: COC(=O)CCC(C)C1CCC2C3C(O[Si](C)(C)C)C=C4CC(O[Si](C)(C)C)CCC4(C)C3CC(O[S
Mol. weight [g/mol]: 637.13

Physical Properties

Property code	Value	Unit	Source
log10ws	-2.41		Crippen Method
logp	9.035		Crippen Method
rinpol	3232.00		NIST Webbook

Sources

Crippen Method: <http://pubs.acs.org/doi/abs/10.1021/ci990307l>
Crippen Method: https://www.chemeo.com/doc/models/crippen_log10ws
NIST Webbook: <http://webbook.nist.gov/cgi/cbook.cgi?ID=R390022&Units=SI>

Legend

log10ws: Log10 of Water solubility in mol/l
logp: Octanol/Water partition coefficient
rinpol: Non-polar retention indices

Latest version available from:

<https://www.chemeo.com/cid/69-598-3/3-7-12-Trihydroxy-5-cholenoic-acid-methyl-ester-TMS.pdf>

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